

SESOIL Execution Using the Data
Management Supporting Systems
SESIN and SESOILFM. User's Guide

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USER'S GUIDE TO
SESOIL EXECUTION USING THE DATA MANAGEMENT
SUPPORTING SYSTEMS, SESIN AND SESOILFM

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I. Introduction

SESOIL is a Seasonal SOIL compartment model developed by Bonazountas and Wagner (1981) for the Office of Pesticides and Toxic Substances (OPTS) of the Environmental Protection Agency (EPA). It estimates the rate of chemical transport/transformation from the soil surface to groundwater (i.e., the unsaturated zone) in terms of mass and concentration distributions among the soil, water (i.e., the unsaturated zone) and air phases. As it was originally designed, SESOIL is batch oriented and, as such, does not support direct user interaction. The preparation and update of input data files can be manually made through the conventional text editor. The output results are in report form containing numerous parameters presented in a fixed-format fashion.

Under the auspices of the modeling team in the Exposure Evaluation Division (EED) of OPTS, the General Software Corporation (GSC) has implemented two data management systems, one for the input and one for the output, in support of the SESOIL modeling operation. The input data management system (SESIN) uses interactive prompts to assist you, the user, in the preparation of the model input files. A set of SESIN commands have been designed to perform efficient file manipulation and data editing. The output data management system (SESILFM) provides for analyses of SESOIL output parameters, using available facilities for graphical, statistical, and tabular displays.

Additionally, a data base is available through SESIN to provide a majority of input data required by SESOIL for a total of seventy (70) cities in the United States. SESILFM can also be used to generate an input data file for AT123D (Yeh, 1981), a transport model for the saturated zone, so that SESOIL and AT123D may be coupled to serve as a complete modeling tool for the subsurface pollutant fate analysis.

The purpose of this document is to describe in detail the procedures required to execute SESOIL through the use of the two data management systems. It is intended as a comprehensive guide for users who have no specialized computer knowledge. However, a working knowledge of the Graphical Exposure Modeling System (GEMS) (GSC 1982), through which SESOIL is executed, is necessary for making more efficient use of the SESOIL model. Only a brief description of SESOIL's input data file structure is presented to facilitate subsequent discussions. More detailed model information is contained in the original SESOIL manual (Bonazountas and Wagner 1981).

II. SESOIL DATA STRUCTURE

A. Input

Three input data files are required by SESOIL: GENERAL, LEVEL, and EXEC. Since the EXEC data file originally required by SESOIL is generated automatically by SESIN, only the GENERAL and LEVEL data files are discussed.

On the basis of the analysis need and the degree of accuracy expected in the modeling results, SESOIL allows you to select among four (4) different types (i.e., LEVEL0, LEVEL1, LEVEL2, AND LEVEL3) of LEVEL data file, each type representing a different level of data requirement sophistication. In general, the LEVEL0 and LEVEL1 data files are for analyses on an annual basis (one year only) of a 2-layer soil column. The LEVEL2 data file is for analyses on a monthly basis (from 1 through 10 years) of a 2-layer soil column. The LEVEL3 data file is similar to the LEVEL2 data file with the exception that the soil column is sliced into a 3-layer configuration. A detailed presentation of these data files is given below.

1. GENERAL DATA FILE

This data file contains the following climatologic, soil, and chemical input parameters:

-- Climatologic

L	Latitude of area (degree N)
TA	Temperature of area (degree C)
NN	Fraction of area covered by clouds (fraction)
S	Relative humidity of area (fraction)
A	Shortwave albedo of surface (-)
REP	Daily evapotranspiration (cm/day)
MPA	Annual precipitation (cm)
MPM	Monthly precipitation (cm)
MTR	Mean time of rain (days)
MN	Mean number of storm events (#)
MT	Mean length of rainy season (days)

(Note: All parameters are presented either yearly or monthly, depending on the selected type of the LEVEL data file. MPA is used for LEVEL0 and LEVEL1 only)

-- Soil

RS	Soil density (g/cm ³)
K1	Soil intrinsic permeability (cm ²)
C	Soil disconnectedness index (-)
N	Effective soil porosity (-)
OC	Organic content of soil (% OC)
CC	Clay content of soil (% CC)
CEC	Soil cation exchange capacity (me/100g)
K1U	Intrinsic permeability of upper soil layer (cm ²)

K1M	Intrinsic permeability of middle soil layer (cm ²)
K1L	Intrinsic permeability of lower soil layer (cm ²)

(Note: All the above parameters are not necessarily needed for a given type of LEVEL data file.)

-- Chemical

SL	Compound solubility in water (ug/ml)
KOC	Adsorption coeff in organic carbon (ug/goc)/(ug/ml)
DA	Diffusion coeff in air (cm ² /sec)
KDE	Biodegradation rate of compound (day ⁻¹)
H	Henry's law constant (m ³ *atm/degK*mol)
K	Adsorption coeff in soil (ug/g)/(ug/m)
MWT	Molecular weight (g/mol)
VAL	Valence
KNH	Neutral hydrolysis constant (day ⁻¹)
KBH	Base hydrolysis constant (L/mol*day)
KAH	Acid hydrolysis constant (L/mol*day)
SK	Stability constant compound-lig and complex(-)
B	Moles of lig and per mole of compound complexed [#]
MWTLIG	Molecular weight of lig and (g/mol)

2. LEVEL0 Data File

This data file contains the following input parameters:

AR	Surface area of the compartment (cm ²)
Z	Depth to groundwater table (m)
DU	Depth of upper soil zone (cm)
PH	pH of upper soil zone (-)
APH	Ratio: pH lower to upper soil zone (-)
AKDE	Ratio: biodegradation rate lower to upper soil zone (-)
AOC	Ratio: Organic carbon lower to upper soil zone (-)
ACC	Ratio: Clay content lower to upper soil zone (-)
ISRA	Index of surface runoff participation in pollutant transport (-)
ASL	Ratio: Pollutant conc. in rain to maximum solubility in water (-)
ACEC	Ratio: CEC lower to upper soil zone (-)
POLINU	Total pollutant load per year entering upper soil zone (ug/cm ²)

POLINL	Total pollutant load per year entering lower soil zone (ug/cm^2)
LIGU	Lig and input mass to upper soil zone (ug/cm^2)
LIGL	Lig and input mass to lower soil zone (ug/cm^2)
THA	Soil moisture content (%)
IA	Infiltration (cm)
RGA	Groundwater recharge (cm)
RSA	Surface runoff (cm)

3. LEVEL1 Data File

This data file contains all the input parameters listed for the LEVEL0 data file except THA, IA, RGA, and RSA.

4. LEVEL2 Data File

This data file contains the following input parameters:

AR	Surface area of the compartment (cm^2)
Z	Depth to groundwater table (m)
DU	Depth of upper soil zone (cm)
PH	pH of upper soil zone (-)
APH	Ratio: pH lower to upper soil zone (-)
AKDE	Ratio: biodegradation rate lower to upper soil zone (-)
AOC	Ratio: Organic carbon lower to upper soil zone (-)
ACC	Ratio: Clay content lower to upper soil zone (-)
FRN	Freundlich equation exponent (-)
ACEC	Ratio: CEC lower to upper soil zone (-)
CUM	Conc. pollutant in soil moisture upper zone (g/ml)
CLM	Conc. pollutant in soil moisture lower zone (g/ml)
POLINU	Pollutant load per unit area entering upper zone (ug/cm^2)
POLINL	Pollutant load per unit area entering lower zone (ug/cm^2)
ISRM	Monthly index for pollutant appearance in surface runoff (-)
ASL	Monthly ratio: conc. of pollutant to maximum solubility in H_2O (-)
TRANSU	Pollutant transformed in upper zone (ug/cm^2)
TRANSL	Pollutant transformed in lower zone (ug/cm^2)
SINKU	Pollutant lost by processes not described (upper) (ug/cm^2)

SINSL	Pollutant lost by processes not described (lower) (ug/cm ²)
LIGU	Lig and mass input to upper zone (ug/cm ²)
LIGL	Lig and mass input to lower zone (ug/cm ²)

(Note: Parameters CUM through LIGL are expressed in terms of monthly values)

5. LEVEL3 Data File

This data file is, in structure, almost identical to LEVEL2 data file with the exception of additional entries for the third (middle) soil layer.

Examples of the GENERAL, LEVEL0, LEVEL1, LEVEL2, and LEVEL3 data files are given in Appendix 3. Note that:

- o The GENERAL data file contains data for different sites, soil types, and chemical compounds.
- o For simulation period of more than one year (LEVEL2 or LEVEL3 application), the input parameters can be repeated for each additional year under consideration for up to 10 years.
- o For simulation of many years of data (up to 100 years) for a desired scenario of pollutant loadings, an auxiliary file (AUX) can be created which contains the variables POLINU, POLINL (LEVEL2) and POLINM (LEVEL3). The user is prompted to build this file only if the number of years of available data in the GENERAL and LEVEL files are less than the desired number of years to be simulated as given by the parameter JYRS in the EXEC data file.
- o LEVEL0, LEVEL1, and LEVEL2 data files also contain data for more than one site location.

B. Output

Output results of SESOIL include hydrologic balance calculations and the distribution of pollutant in the soil column. After successfully following the procedures described in this document, you are able to obtain a complete hard-copy SESOIL output for review. A total of 86 selected SESOIL output parameters can be stored in GEMS for further analysis.

III. CONCEPT OF DEFAULT FILES

To minimize the complications involved in the preparation of SESOIL input data files, the design of SESIN is based on the central concept of providing you with a complete set of default files with which you can perform all necessary file editing and manipulations. In requesting the default files, there are three options supported by SESIN:

1. The default files are retrieved from the 70-city data base (Note: This data base only provides data for Level 3 operation).
2. The default files are retrieved from the files presented in Appendix 3.
3. The default files are specified by you.

The use of Option 1 is illustrated in Appendix 1. Note that once these default files are available to you, you must review all the default values carefully to check their applicability to your simulation run. In some cases, some default values are provided for you only to ensure the completeness of the file content and they must be revised by you in order to give meaningful simulation results.

IV. GUIDE TO SESOIL EXECUTION AND DATA MANAGEMENT

A. Access to SESOIL

You can access SESOIL through the Graphical Exposure Modeling System (GEMS). The host computer of GEMS is Digital Equipment Corporation VAX-11/780, located at EPA's National Computer Center (NCC) in Research Triangle Park, North Carolina.

It is assumed that you are familiar with the log-on procedures of the VAX system and are able to invoke GEMS (Consult User's Guide to GEMS, if not). After you make connection, the two steps required to gain access to SESOIL in GEMS are:

ENTER AN OPERATION: . M <CR>

ENTER MODEL NAME: SESOIL <CR>

- * (Where <CR> stands for pressing the carriage return key on the terminal).

B. Major Decision Points in SESOIL Execution

At the outset, and through the process of SESOIL execution, you are confronted with the following major decision points:

- (1) Do I have all the GENERAL, LEVEL (LEVEL0, LEVEL1, LEVEL2, or LEVEL3), EXEC, and AUX (if necessary) data files ready for execution, so that I do not need to create those data files through SESIN? (Note: (a) The EXEC data file is required if SESIN is not used. Refer to the original SESOIL manual for description. (b) The AUX file is required if JYRS in the EXEC data file is greater than the number of years of data available in the LEVEL and GENERAL data files.)
- (2) If I use SESIN support, which of the three options (see Section III) do I choose to get my default GENERAL and LEVEL data files?

- (3) Do I want to generate additional years of pollutant loading data?
- (4) Do I want to make multiple SESOIL runs?
- (5) Do I want the support of the output data management system, SESOILFM, so that after the SESOIL execution is completed, I can perform graphical, statistical, and tabular output manipulations using GEMS' facilities?
- (6) If the answer to (5) is "yes", what is the dataset name (1 to 12 characters) under which I want my SESOIL output to be stored in GEMS?
- (7) Do I want to request an AT123D input file based on the SESOIL simulation results?
- (8) If the answer to (7) is "yes", do I want to make an AT123D run using the AT123D input file just created?

A flowchart outlining the relative sequence of the above decision points is presented in Figure 1. As shown, the entire process can be divided into five execution branches which will be discussed below.

C. Detailed Guide to SESOIL Execution

To facilitate the following discussion, this section will be presented in terms of the five execution branches shown in Figure 1.

1. Branch 1

This is the core of the GSC-developed SESOIL data management system. The execution will enter this branch if you choose to create new data files (i.e. GENERAL and LEVEL data files) through SESIN. The main idea is to guide you interactively through an efficient data editing and file preparation process. To do this, SESIN is equipped with the following user commands:

ALL	BRIEF	<CR>	DONE	EXIT	HELP
LA120	LAST	MENU	NEXT	NOBRIEF	NONE
REST	REVIEW	REVISE	VT100		

A detailed list of command descriptions is presented in Table 1.

cities

To start with, you need to select one of the three options for getting you default data files (see Section III). After the selection, you are guided by self-explanatory prompts provided by SESIN. You can respond with brief answers or use the above listed commands to control the progress of the process. The use of SESIN and its commands can be illustrated best by Appendix 1, which contains a step-by-step example side-marked with comments. You are advised to follow this example carefully in order to make efficient use of SESIN. It is to be stressed that Appendix 1 only

presents an entry-level demonstration of the use of SESIN. After becoming more familiar with the organization of SESOIL input files and the concepts of SESIN, you are encouraged to explore the other powerful SESIN capabilities through hands-on training practices.

1.a. Branch 1-A

This branch is a sub branch of Branch 1. It represents the use of generating long-term pollutant application scenarios. The execution will enter this branch if the value of JYRS (desired number of years to be simulated) in the EXEC data file (as specified by the user) is greater than the number of years of data available in the GENERAL and LEVEL data files. If this is true, the user is prompted to specify the additional pollutant loading data (POLINU, POLINM, POLINL) for the additional years required, which is then stored in the auxiliary data file AUX. The remaining parameter values of the GENERAL and LEVEL data files for the additional years are automatically generated by repeated use of the last year of data. An example run using this Branch is given in Appendix 2.

2. BRANCH 2

This branch represents the original SESOIL data file input fashion in that it prompts you to input directly the three data files: GENERAL, LEVEL (i.e., LEVEL0, LEVEL1, LEVEL2, or LEVEL3), and EXEC. The auxiliary pollutant loading data file, AUX is also prompted for. Note that the user-prepared EXEC data file is needed in this branch. The SESOIL execution will enter this branch if you select not to create new data files through SESIN. An example is given in Appendix 3 to illustrate the use of this branch.

3. BRANCH 3

This branch represents the use of the GSC-developed SESOIL output data management system (SESOILFM), which allows you to create a dataset containing 86 selected SESOIL output variables, as shown in Table 2. This dataset will be stored in GEMS so that you can utilize GEMS' capabilities for output data manipulation, analysis, and interpretation. The columns of VAR NAME and VAR UNIT in Table 2 give brief variable definitions and variable units, respectively, to assist you in recognizing the variables to be selected for further review. Note that Variables 5 through 7, representing chemical half-life periods for different soil layers are not direct SESOIL output variables. Instead, they are derived from the degradation rates (Variables 2 through 4) of chemicals assuming a first order decay. Furthermore, Variables 73 through 86, each representing a possible transformation/transport process in the pollutant fate analysis, are calculated from the SESOIL output to show the total mass balance distribution in the unstaured soil zone at the end of each year being simulated. Each of the variables in Table 2 is stored in the dataset on a monthly basis (i.e., not monthly applicable, such as PRECIP, MASGRW, etc.), their estimated values are stored in the dataset at the 12th month of each year. Spaces prior to the 12th month are flagged as "NO VALUE".

In addition to the variables shown in Table 2, the following identification and time-factor variables are included in the dataset to facilitate the use of GEMS' capabilities.

RUN:	Run identification number
COMP ID:	Compound name studied
LEVEL:	Modeling level selected
YEAR:	Year simulated
MONTH:	Cumulative month through the years simulated
MINDEX:	Month of given year (non-cumulative)
VAR ID:	Variable identification code (to be used to call out the variable needed)
VAR NAME:	More descriptive variable definition
VAR UNIT:	Unit for each variable

As shown, each of these variables serves a specific identification purpose. You can also use the variable RUN to retrieve (via GEMS EXTRACT and/or FILTER function) the variables for a particular run from a dataset containing multiple run records. The variable MONTH provides an across-the-years monthly time index, which you can use to observe the continuous variation of a given variable throughout the years being analyzed.

There are only three steps required to use SESOILFM for creating a SESOIL output dataset in GEMS, as exemplified near the end of Appendix 1. The steps are:

(i) enter Y in response to the prompt: "Enter Y to output SESOIL results to GEMS>"

(ii) give a dataset name (1 to 12 characters) in response to the prompt: "Enter GEMS output dataset name>"

(iii) enter a dataset description (1 to 60 characters) in response to the prompt: "Enter tag field>."

Any characters other than Y entered in Step (i) cause the execution to skip the the SESOILFM supporting service. The tag field is optional and you use it only for the purpose of more descriptive dataset identification. You can skip this entry by pressing <CR>.

It is beyond the scope of this document to describe the applications of GEMS' graphical, statistical, and tabular facilities to the SESOIL dataset described above. You are advised to consult the User's Guide to GEMS (GSC 1982) for detailed information. However, to demonstrate what kind of products you may obtain from the use of SESOILFM and GEMS file management facilities, the following examples are presented:

Figure 2: This is a display of the predicted monthly pollutant mass transport into the groundwater (INMOGRW) vs. the cumulative month index (MONTH) for a period of four years. The two curves, presenting a vivid visual comparison, are derived from two separate SESOIL runs, one with a biodegradation rate (KDE) of 1.0 day^{-2} and the other of 0.22 day^{-2} .

Figure 3: This is a bar chart showing the cumulative effect of the curve with a KDE of 1.0 day^{-2} in Figure 2.

Table 3: This is a GEMS' tabular output presenting the four variables: MONTH, MINDEX, CONWMOU, MASWMOU. Note that Table 2 is also produced using SESOILFM and GEMS.

4. BRANCH 4

This branch is activated if you elect to bypass SESOILFM by entering any characters other than Y in response to the prompt: "Enter Y to output SESOIL results to GEMS>". An example of this is shown near the end of Appendix 1.

5. BRANCH 5

This branch allows you to review the SESOIL results that are applicable to the construction of an AT123D input file. After review, if you are satisfied with the results, you can request the build up of an AT123D input file. Due to the complexity of the file content, the following assumptions are used:

- o The pollutant source area is assumed to be a square with a size equivalent to the soil column area used in SESOIL.
- o The aquifer depth and width are assumed to be infinite.
- o The hydraulic gradient is assumed to be 0.01.
- o The hydraulic conductivity is based on SESOIL's intrinsic permeability.
- o The longitudinal, transverse, and vertical dispersivities are assigned on the basis of Yeh's (1981) typical values and the calculated hydraulic conductivity.
- o The time interval size is assumed to be 8760 hours (i.e., one year for Level 2 and 3; 730 hours (i.e., one month) for level 0 and 1.

- o The loading source release is assumed to be constant (NSOUS=0) for level 0 and 1; time-dependent (NSOUS=1) for level 2 and 3.
- o The loading source release is assumed to be instantaneous (INSTAN=0) for level 0 and 1; continuous (INSTAN=1) for level 2 and 3.
- o The time duration of waste release is assumed to continue through the entire SESOIL simulation period for level 2 and 3; not used for level 0 and 1.
- o The variable source release rate, calculated by SESOIL, is used for level 2 and 3.
- o The decay constant is set equivalent to the neutral hydrolysis rate used in SESOIL.
- o The distribution coefficient is taken as equal to the adsorption coefficient used in SESOIL.

Several of the assumptions which may require special attention are printed as "Notes" in the Branch-5 output as shown in the sample run in Appendix 1.

In addition to the above assumptions, the AT123D input file is also developed using various assumed program functional and execution control values. You must carefully review the file content to ensure the applicability of all input values to your specific need. You can edit or execute this AT123D input file using the AT123D data management systems, AT123DIN and AT123DOUT, by responding to the last prompt in Appendix 1 with a "yes". For more information concerning the use of AT123D data management systems, consult the AT123D User's Guide (GSC, 1983).

V. SUMMARY

A concise guide to SESOIL execution using GSC-developed input (SESIN) and output (SESOILFM) data management systems is presented in this document. The basic SESOIL data structure is discussed briefly to provide the necessary background for the efficient use of SESIN and SESOILFM. The execution logic is outlined by five major branches, each of which may be invoked or bypassed by you using interactive commands. Most importantly, detailed examples (Appendix 1 and Appendix 2) with step-by-step explanatory comments are provided to guide the user through a successful test SESOIL execution. The user is advised to follow those examples closely to ensure the most efficient use of the systems.

Again, it is to be stressed that Appendices 1, 2 and 3 present only an entry-level demonstration of the use of SESIN and SESOILFM, respectively. The more advanced capabilities of these two systems can be learned via hands-on trial practices. For assistance, call GSC at (301) 459-9494.

VI. ACKNOWLEDGEMENTS

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VII. REFERENCES

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TABLE 1. SESIN COMMAND SUMMARY

The following is a list of commands that apply within SESIN. Note that all commands do not work at all prompts.

ALL	BRIEF	<CR>	DONE	EXIT	HELP	LA120
LAST	MENU	NEXT	NOBRIEF	NONE	REST	REVIEW
REVISE	VT100					

for additional help with a command type: HELP <COMMAND>

for a full list of commands with explanations type: HELP *

- | | |
|-------|---|
| ALL | - for parameter input ALL designates that all parameters will be changed.
- in response to a request for a list of months ALL designates that values for all 12 months will be set to a single input value. |
| BRIEF | - no parameter descriptions will be given when a parameter menu is displayed.
- a carriage return assumes the currently selected data. |
| DONE | - previous input and existing values are accepted and program execution continues with the next prompt. |
| EXIT | - the program run is immediately ended if no runs have yet been fully set up. If at least one run has been set up you will go through the exit procedure which will create your run files. If you type exit again at this point the program run will be immediately stored. |
| HELP | - provides command summary and descriptions. |
| LA120 | - informs the program that the user is on an LA120 or other hardcopy terminal. Program output is slightly modified for this terminal type. This is the default terminal type. |
| LAST | - this command can be used to modify climatic or level data for the year preceeding the current year. This is particularly useful when using the revise command to modify data from a prior run. |

TABLE 1. SESIN COMMAND SUMMARY(CONTINUED)

MENU	- displays the current parameter list when at a data item prompt.
NEXT	- this command is the complement to the LAST command and can be used to move forward a year in a climatic or level data set.
NOBRIEF	- allows display of parameter descriptions to be restored after BRIEF has been used.
NONE	- at site, chemical, or soil selection this command allows the zeroing of the input data array for entry of an entirely new set of data. - at all other prompts NONE is identical to DONE
REST	- allows the user to copy the remainder of the default input data file to the sesoil input file being developed. This command only works at the enter data item to be updated prompt for monthly level or climatic data.
REVIEW	- allows the user to review the current data set. The user is prompted to enter C or A for review to review either data for just the current year or for all years up to the current year.
REVISE	- allows modification of the currently selected data (level, climatic, soil or chemical). This command is used when making multiple runs which require modification of data selected in the previous run. - NOTE when this command is used when in level 2 or 3 (monthly data) you will be left at the last data year entered when altering monthly parameters.
VT100	- this command turns off the LA120 command and can be used for all CRT terminals.

TABLE 2. SESOILFM VARIABLES SUMMARY

VAR ID	VAR NAME	VAR UNIT	VAR ID	VAR NAME	VAR UNIT
1	DEPTHGRW		51	MASWRL	
2	DEGRDU	DEPTH TO GRW	52	MASWRL	YR MASS-MOIS(L)
3	DEGRDM	DEGRAD. RATE(U)	53	MASWRL	MO MASS-MOIS(U)
4	DEGRDL	DEGRAD. RATE(M)	54	MASWRL	MO MASS-MOIS(M)
5	HALFU	DEGRAD. RATE(L)	55	MASWRL	YR MASS-MOIS(L)
6	HALFM	HALF-LIFE(U)	56	MASWRL	YR MASS-MOIS(M)
7	HALFL	HALF-LIFE(M)	57	MASWRL	YR MASS-SOIL(U)
8	HALFL	HALF-LIFE(L)	58	MASWRL	YR MASS-SOIL(M)
9	ADSOIL	ADSORP CO(SOIL)	59	MASWRL	MO MASS-SOIL(L)
10	ZOCU	ADSORP CO(O.C.)	60	MASWRL	MO MASS-SOIL(M)
11	ZOCM	% ORG. CARB.(U)	61	MASWRL	MO MASS-SOIL(L)
12	ZOCL	% ORG. CARB.(M)	62	MASWRL	YR CONC-AIR(U)
13	HENRY	% ORG. CARB.(L)	63	MASWRL	YR CONC-AIR(M)
14	DENSITY	HENRY CONSTANT	64	MASWRL	MO CONC-AIR(L)
15	PERMALL	SOIL DENSITY	65	MASWRL	MO CONC-AIR(M)
16	PERMU	INTRA PERM(AVE)	66	MASWRL	YR CONC-AIR(L)
17	PERMH	INTRA PERM(U)	67	MASWRL	YR CONC-AIR(M)
18	PERML	INTRA PERM(M)	68	MASWRL	YR MASS-AIR(U)
19	PEROSITY	INTRA PERM(L)	69	MASWRL	YR MASS-AIR(M)
20	PRECIP	SOIL POROSITY	70	MASWRL	MO MASS-AIR(L)
21	INRYU	YEARLY PRECIP.	71	MASWRL	MO MASS-AIR(M)
22	INRYM	POL YR INPUT(U)	72	MASWRL	MO MASS-AIR(L)
23	INRYM	POL MO INPUT(U)	73	MASWRL	MO MASS-AIR(M)
24	INRYM	POL YR INPUT(M)	74	MASWRL	MO MASS-AIR(L)
25	INRYL	POL MO INPUT(M)	75	MASWRL	MO MASS-AIR(M)
26	INRYL	POL YR INPUT(L)	76	MASWRL	MO MASS-AIR(L)
27	MOIST	POL MO INPUT(L)	77	MASWRL	MO MASS-AIR(M)
28	GRUFLOW	SOIL MOISTURE	78	MASWRL	MO MASS-AIR(L)
29	DEGRYU	GRW RUNOFF	79	MASWRL	MO MASS-AIR(M)
30	DEGRYU	YR DEGRADED(U)	80	MASWRL	MO MASS-AIR(L)
31	DEGRYU	MO DEGRADED(U)	81	MASWRL	MO MASS-AIR(M)
32	DEGRYU	YR DEGRADED(M)	82	MASWRL	MO MASS-AIR(L)
33	DEGRYU	MO DEGRADED(M)	83	MASWRL	MO MASS-AIR(M)
34	DEGRYU	YR DEGRADED(L)	84	MASWRL	MO MASS-AIR(L)
35	INRYGRW	MO DEGRADED(L)	85	MASWRL	MO MASS-AIR(M)
36	INRYGRW	YR INTO GRW	86	MASWRL	MO MASS-AIR(L)
37	CONWYU	MO INTO GRW			
38	CONWYU	YR CONC-MOIS(U)			
39	CONWYU	YR CONC-MOIS(M)			
40	CONWYU	YR CONC-MOIS(L)			
41	CONWYU	MO CONC-MOIS(U)			
42	CONWYU	MO CONC-MOIS(M)			
43	CONWYU	MO CONC-MOIS(L)			
44	CONWYU	YR CONC-SOIL(U)			
45	CONWYU	YR CONC-SOIL(M)			
46	CONWYU	YR CONC-SOIL(L)			
47	CONWYU	MO CONC-SOIL(U)			
48	CONWYU	MO CONC-SOIL(M)			
49	CONWYU	MO CONC-SOIL(L)			
50	MASWRL	YR MASS-MOIS(U)			
		YR MASS-MOIS(M)			
		YR MASS-MOIS(L)			

U: UPPER SOIL LAYER

M: MIDDLE SOIL LAYER

L: LOWER SOIL LAYER

YR: YEARLY

MO: MONTHLY

TABLE 3. EXAMPLE OF SESOILFM PRODUCT(TABULAR DISPLAY)

#	MONTH	MINDEX	CONWMOU	MASWMOU
1	1	1	0.37293001E-02	86906.555
2	2	2	0.74580982E-02	173801.42
3	3	3	0.11186397E-01	260684.66
4	4	4	0.14914199E-01	347556.31
5	5	5	0.18641606E-01	434418.75
6	6	6	0.22368507E-01	521269.41
7	7	7	0.26095007E-01	608110.75
8	8	8	0.29821001E-01	694940.19
9	9	9	0.33546392E-01	781755.69
10	10	10	0.37271384E-01	868561.88
11	11	11	0.40995900E-01	955356.94
12	12	12	0.44719987E-01	1042142.0
13	13	1	0.48443485E-01	1128913.4
14	14	2	0.52166484E-01	1215673.1
15	15	3	0.55889081E-01	1302423.4
16	16	4	0.59611194E-01	1389162.5
17	17	5	0.63332774E-01	1475889.1
18	18	6	0.67053966E-01	1562606.8
19	19	7	0.70774674E-01	1649313.1
20	20	8	0.74494869E-01	1736007.4
21	21	9	0.78214593E-01	1822690.9
22	22	10	0.81933878E-01	1909364.1
23	23	11	0.85652687E-01	1996026.1
24	24	12	0.89371011E-01	2082676.9
25	25	1	0.93088754E-01	2169314.0
26	26	2	0.96806012E-01	2255940.0
27	27	3	0.10052200	2342536.5
28	28	4	0.10423800	2429133.0
29	29	5	0.10795202	2515683.5
30	30	6	0.11166400	2602186.5
31	31	7	0.11537600	2688689.8
32	32	8	0.11908799	2775193.3
33	33	9	0.12280001	2861697.0
34	34	10	0.12651201	2948200.3
35	35	11	0.13022399	3034703.3
36	36	12	0.13393600	3121207.0

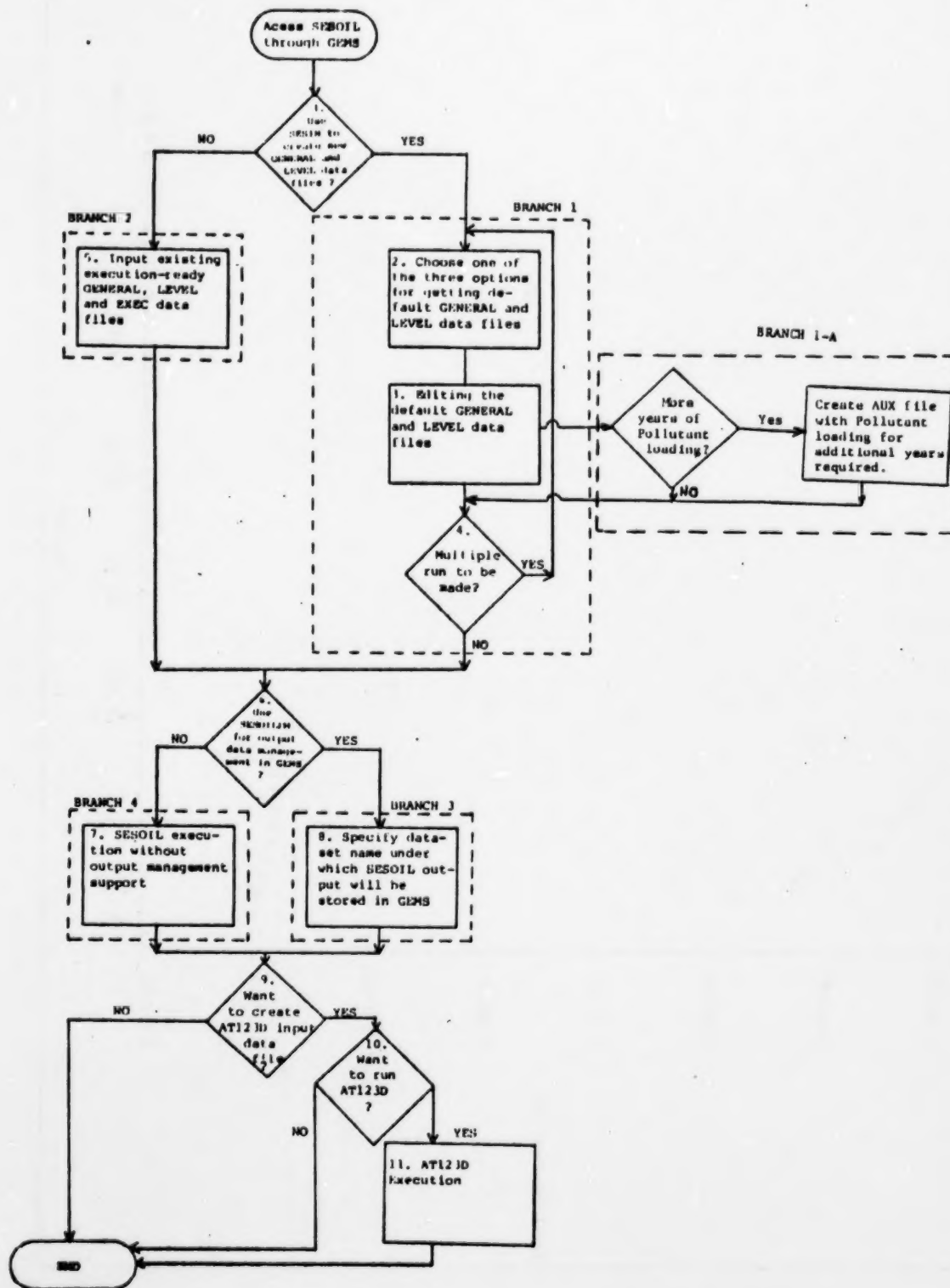


FIGURE 1. FLOWCHART FOR SESOIL EXECUTION WITH DATA MANAGEMENT SYSTEM SUPPORT

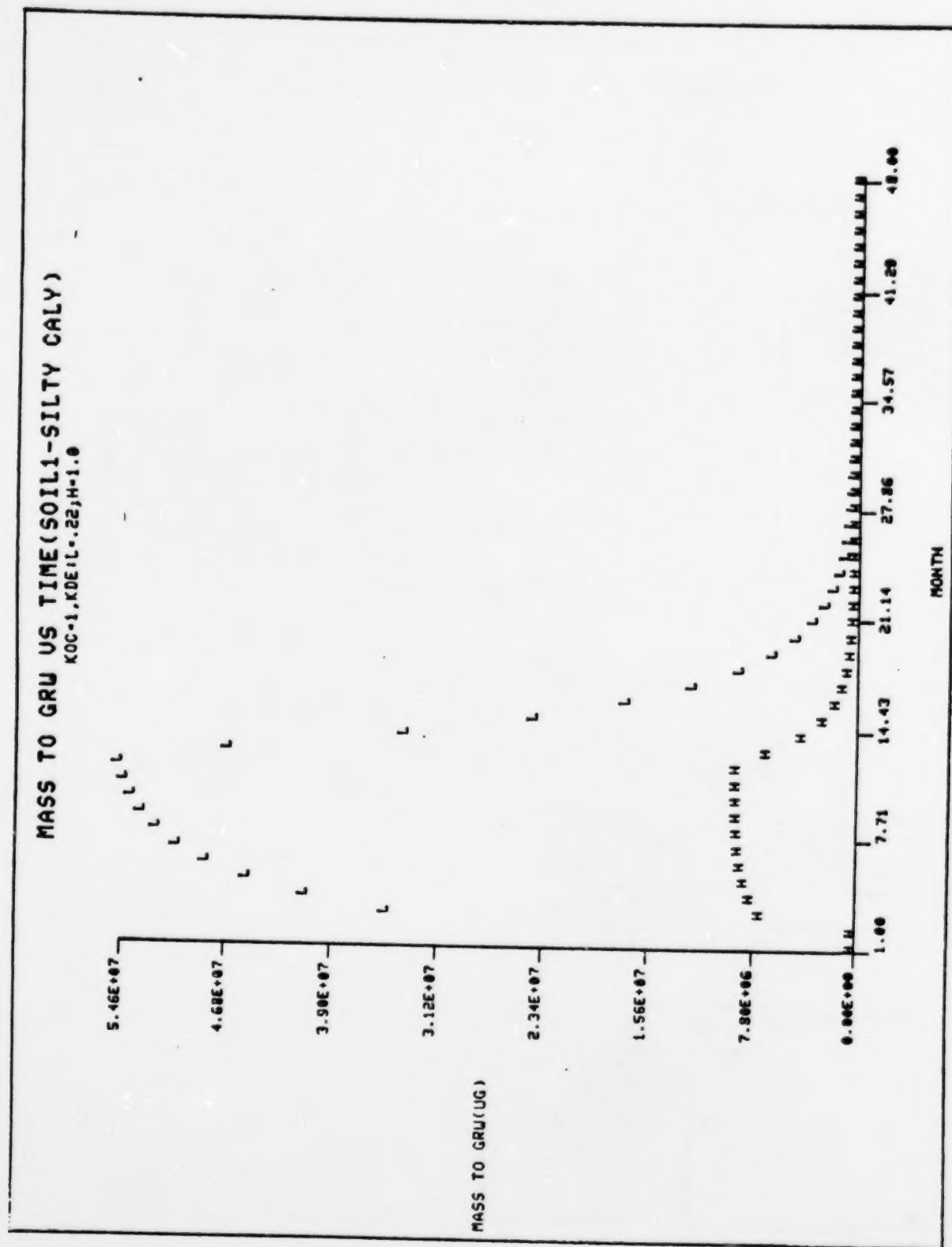


FIGURE 2. EXAMPLE OF SESOILF PRODUCT (SCATTER DIAGRAM)

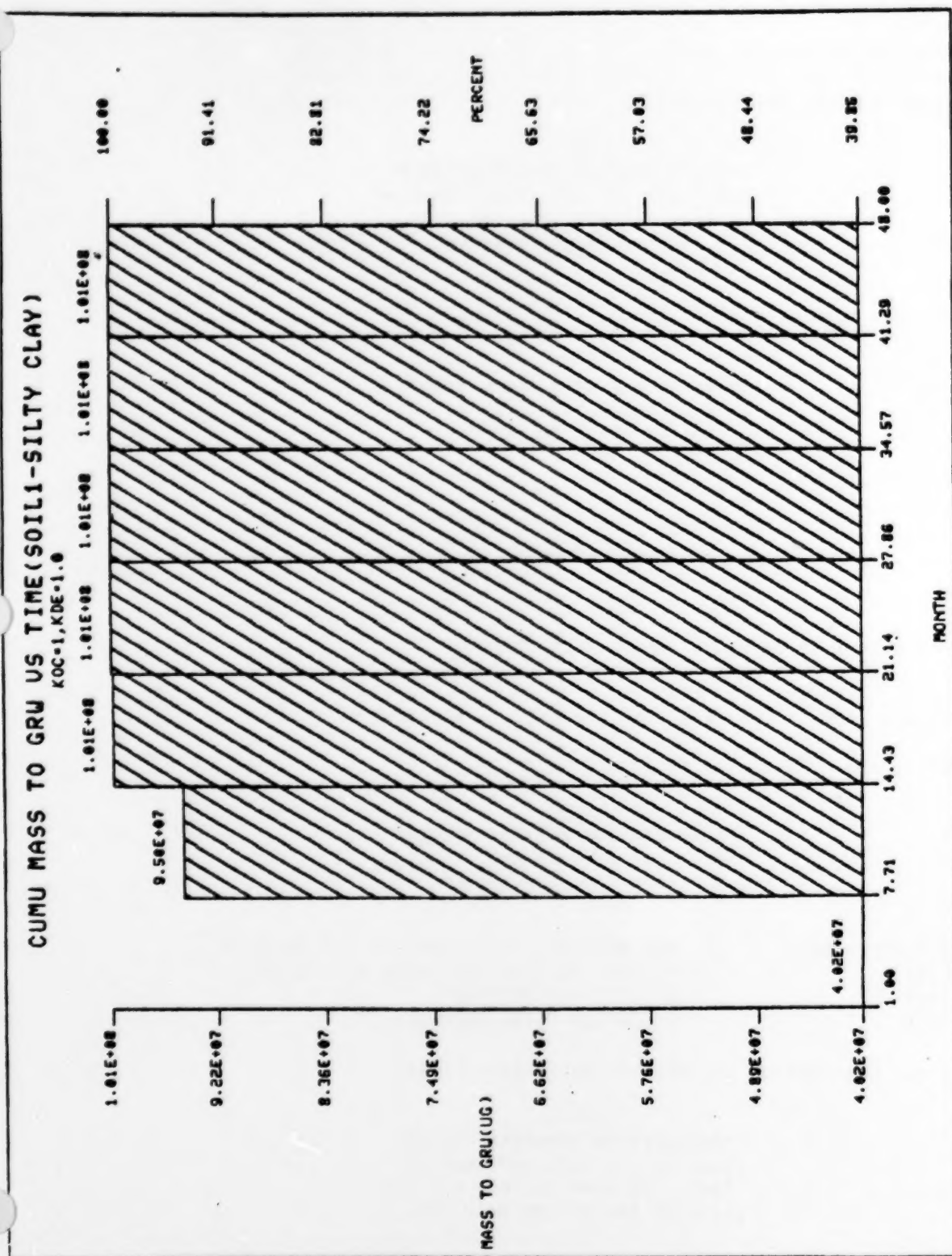


FIGURE 3. EXAMPLE OF SESOILFM PRODUCT(CUMULATIVE BARCHART)

APPENDIX 1: SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM

Enter an operation: [m]

Enter a model name: [sesoil]

NOTE

These are the two steps required
to access SESOIL through GEMS.
[] encloses your response.

Are you creating SESOIL input files: [yes]

NOTE

This is the decision to enter
either BRANCH 1 or BRANCH 2
(see figure 1).

SEIN
INTERACTIVE SESOIL INPUT
GSC VER1.1
5-DEC-84

For a list of commands type : HELP COMMANDS

Do you want to use data from the cities database? [yes]

Enter the name of the desired data site: [help]

NOTE

'help' was used to get
assistance from the system.

A PLACE NAME	THE NAME OF A DATA SITE IN THE DATABASE
LIST	TO LIST ALL THE AVAILABLE DATA SITES
EXIT	TO EXIT THE PROGRAM
HELP	TO OBTAIN THESE MESSAGES

Enter the name of the desired data site: [list]

NOTE

Based on the help information,
'list' was used to show all the
sites in the cities database.

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

Enter the state abbreviation or ALL for entire list [all]

NOTE

A list containing all sites in
all states was requested.

DATA AVAILABLE FOR:

MOBILE AL	JUNEAU AK	PHOENIX AZ
LITTLE ROCK AR	LOS ANGELES CA	SACRAMENTO CA
SAN FRANCISCO CA	DENVER CO	HARTFORD CT
WILMINGTON DE	WASHINGTON DC	JACKSONVILLE FL
MIAMI FL	ATLANTA GA	HONOLULU HI
BOISE ID	CHICAGO IL	PEORIA IL
INDIANAPOLIS IN	DES MOINES IA	WICHITA KS
LOUISVILLE KY	NEW ORLEANS LA	PORTLAND ME
BALTIMORE MD	BOSTON MA	DETROIT MI
SAULT STE MARIE MI	DULUTH MN	MINNEAPOLIS/ST PAUL MN
JACKSON MS	KANSAS CITY MO	ST LOUIS MO
GREAT FALLS MT	OMAHA NB	RENO NV
CONCORD NH	ATLANTIC CITY NJ	ALBUQUERQUE NM
ALBANY NY	BUFFALO NY	NEW YORK NY
CHARLOTTE NC	RALEIGH NC	BISMARCK ND
CINCINNATI OH	CLEVELAND OH	COLUMBUS OH
OKLAHOMA CITY OK	PORTLAND OR	PHILADELPHIA PA
PITTSBURGH PA	PROVIDENCE RI	COLUMBIA SC
SIOUX FALLS SD	MEMPHIS TN	NASHVILLE TN
DALLAS/FORT WORTH TX	EL PASO TX	HOUSTON TX
SALT LAKE CITY UT	BURLINGTON VT	NORFOLK VA
RICHMOND VA	SEATTLE/TACOMA WA	SPOKANE WA
CHARLESTON WV	MILWAUKEE WI	CHEYENNE WY
SAN JUAN PR		

Enter the name of the desired data site: [boston]

CURRENT SITE: BOSTON MA

Select run option: [help]

BACK	USE TO RETURN TO THE LIST OF SITES
BUILD	USE TO BUILD SESOIL/TOXSCREEN INPUT DATA FILES
REVIEW	USE TO REVIEW DATA FOR THE SELECTED SITE
EXIT	USE TO EXIT THE PROGRAM

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

CURRENT SITE: BOSTON MA

Select run option: [build]

NOTE

Request was made to build up the
SESOIL input data files from the
70-city database for the Boston
site.

Enter soil data option: [help]

DEFAULT	THE PROGRAM WILL SELECT AN APPROPRIATE SOIL
SPECIFIC	USER SELECTS A SOIL FROM SITE SPECIFIC SOILS
GENERIC	USER SELECTS A SOIL FROM A LIST OF GENERIC SOILS
BACK	TO RETURN TO THE PREVIOUS PROMPT
HELP	TO OBTAIN THESE HELP MESSAGES

Enter soil data option: [default]

NOTE

Request was made to let the
system automatically select an
appropriate soil which is one of
the needed SESOIL input information.

Do you wish to build a SESOIL data set for another location? [no]

NOTE

To simplify the illustration,
only one site (Boston) is
selected here.

Enter your name for the general data file: [gen]

Enter your name for the level data file: [lev]

NOTE

You can name the GENERAL and
LEVEL data files which contain
the data accessed from the
cities database.

** Database access completed **

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFH(CONT'D)

Current GENERAL data file is : GEN
To use an alternate GENERAL data file,
enter the file name (<CR> for none)? [<CR>]

NOTE

An option to let you call in
another GENERAL file to be used
instead of the newly created
file:GEN.DAT, <CR> indicates that
you want to bypass this option.

***** DETERMINATION OF SESOIL POLLUTANT APPLICATION LEVEL *****

City DATABASE access

Selected SESOIL application level is : 3

NOTE

Remember that the 70-city data-
base can be used to build up
files for LEVEL 3 study only.
For other Levels, select the
remaining two options (see Sec.III).

***** AVAILABLE CLIMATIC REGIONS *****

1. BOSTON MA CANTON SANDY LOAM FINE S

Only one site is available ** automatically selected

***** CLIMATOLOGICAL PARAMETERS *****

1. L [des N] latitude of area
2. NAME

Enter data item to be updated : [review]

NOTE

You are strongly recommended to
review all retrieved data to
check their validity.

You may review data for the current year or for all years available.

APP.1:SESOIL~SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

Enter [C]urrent or [A]ll? [all]

NOTE

Since there is only one year of
data presently available, the
entry of 'C' or 'A' makes no
difference.

	1 BOSTON MA	CANTON SANDY LOAM FINE S											1
L	42.36												
TA	13.00	7.330	.5556	-1.56	-0.89	3.39	9.22	14.78	20.00	22.94	21.83	18.06	
NN	0.53	0.6	0.6	0.6	0.56	0.56	0.58	0.58	0.57	0.56	0.53	0.51	
S	0.57	0.61	0.61	0.57	0.57	0.57	0.53	0.57	0.6	0.56	0.57	0.61	
A	0.18	0.21	0.28	0.37	0.35	0.28	0.21	0.18	0.18	0.18	0.18	0.18	
REP	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	
MPH	7.67	11.46	10.77	9.37	8.99	10.19	8.86	8.81	8.10	6.96	8.79	8.03	
MTR	0.6667	1.040	.7083	0.375	0.750	.2917	0.3750	.33330	.20830	.54170	.0833	0.5	
MN	7.00	4.00	9.00	6.00	8.00	5.00	9.00	7.00	8.00	9.00	5.00	7.00	
MT	9.00	11.00	12.00	12.00	11.00	12.00	11.00	11.00	11.00	9.00	10.00	9.00	

Press return to continue? [CR]

Enter data item to be updated : [done]

NOTE

The use of 'done' signals that
the editing of the above data
has been completed and you are
advanced to the next prompt.

Current YEAR : 1

1. TA [deg C] temp of area
2. NN [fraction] fraction of area covered by clouds
3. S [fraction] relative humidity of area
4. A shortwave albedo of surface
5. REP [cm/day] daily evapotranspiration or 0
6. MPH [cm] monthly precipitation
7. MTR [days] mean time of rain
8. MN [#] mean number of storm events
9. MT [days] mean length of rainy season

Enter data item to be updated : [1]

NOTE

'1' means the values for TA are
to be updated.

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

The current values for TA [deg C] are

1 Oct	2 Nov	3 Dec	4 Jan	5 Feb	6 Mar
13.00	7.33	0.55559999	-1.56	-0.89	3.39
7 Apr	8 May	9 June	10 July	11 Aug	12 Sept
9.22	14.78	20.00	22.94	21.83	18.06

Enter the month(s) to alter? [10]

NOTE

The value for July is to be
changed.

Enter the new value(s)? [24.0]

Current YEAR : 1

Enter data item to be updated : [done]

Would you like to enter data for another year? [no]

NOTE

To simplify the illustration,
only one year of data is selected
here. Multiple-year runs are
shown in Appendix 2.

***** AVAILABLE SOILS *****

1. CANTON SANDY LOAM FINE SANDY LOAM BOSTON

Only one soil is available ** automatically selected

***** SOIL PARAMETERS *****

1. RS	[g/cm2]	soil density
2. K1	[cm2]	soil intrinsic permeability
3. C		soil disconnectedness index
4. N		effective soil porosity
5. OC	[%oc]	organic content of soil
6. CC	[%cc]	clay content of soil
7. CEC	[me/100g]	soil cation exchange capacity
8. K1U	[cm2]	intrinsic permeabilities of upper soil layer

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

- 9. K1M [cm2] intrinsic permeabilities of middle soil layer
- 10. K1L [cm2] intrinsic permeabilities of lower soil layer
- 11. RDUST [us/m3] dust loading factor for TOXSCREEN only
- 12. NAME

Enter data item to be updated : [review]

NOTE

Review is made here again to
to check the data.

1 CANTON SANDY LOAM FINE SANDY LOAM BOSTON

- RS,K1,C,N,OC,CC	1.35	0.00	6.00	0.26	0.00	0.00
- CDC,K1U,K1M,K1L,RDUST	0.00.288E-7.288E-7.936E-7	0.00				

Press return to continue? [<CR>]

Enter data item to be updated : [5]

The current value for OC is : 0.00 [%oc]
Enter the new value, or hit <CR> to leave the same : [0.20]

NOTE

The value of OC is changed from
0.00 to 0.20.

Enter data item to be updated : [done]

***** AVAILABLE CHEMICALS *****

1. TEST CHEMICAL

Only one chemical is available ** automatically selected

***** CHEMICAL PARAMETERS *****

- 1. SL [us/mL] compound solubility in water
- 2. KOC [(us/s oc)/(us/mL)] adsorption coeff in organic carbon
- 3. DA [cm2/sec] diffusion coefficient in air
- 4. KDE [day-1] biodegradation rate of compound
- 5. H [m3*atm/desK*mol] Henrys Law constant
- 6. K [(us/s)/(us/mL)] averaged adsorption coeff for compound on soi
- 7. MWT [g/mol] molecular weight
- 8. VAL valence
- 9. KNH [day-1] neutral hydrolysis constant
- 10. KBH [L/mol*day] base hydrolysis constant
- 11. KAH [L/mol*day] acid hydrolysis constant

12. SK stability constant compound-lisand complex
 13. R [#] moles of lisand per mole of bound compound
 14. MWTLIG [g/mol] molecular weight of lisand
 15. NAME

Enter data item to be updated : [review]

1 TEST CHEMICAL

SL,KOC,DA,KDE,HENRY,K	730.00	157.00	0.0825	0.002	0.0031	0.00
MWT,VAL,KNH,KBH,KAH	133.40	0.00	0.00	0.00	0.00	
SK,B,MWTLIG	0.00	0.00	0.00			

Press return to continue? [<CR>]

Enter data item to be updated : [all]

NOTE

 After review, changes to all of
 the above data are requested.
 The changes are carried out item
 by item, as shown below.

NOTE

 Normally you would want to
 replace the TEST CHEMICAL
 parameters with your own
 chemical data.

The current name is : TEST CHEMICAL
 Enter the new name : [pcb-cl2h9cl]

The current value for SL is : 730.00 [ug/mL]
 Enter the new value, or hit <CR> to leave the same : [1.17]

The current value for KOC is : 157.00 [(ug/g oc)/(ug/mL)]
 Enter the new value, or hit <CR> to leave the same : [11500.0]

The current value for DA is : 0.0825 [cm²/sec]
 Enter the new value, or hit <CR> to leave the same : [0.06]

The current value for KDE is : 0.002 [day⁻¹]
 Enter the new value, or hit <CR> to leave the same : [0.0346]

The current value for H is : 0.0031 [m³atm/desK*mol]
 Enter the new value, or hit <CR> to leave the same : [0.0059]

The current value for K is : 0.00 [(ug/d)/(ug/mL)]
 Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for MWT is : 133.40 [g/mol]
Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for VAL is : 0.00
Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for KNH is : 0.00 [day-1]
Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for KBH is : 0.00 [L/mol*day]
Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for KAH is : 0.00 [L/mol*day]
Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for SK is : 0.00
Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for R is : 0.00 [#]
Enter the new value, or hit <CR> to leave the same : [<CR>]

The current value for MWTIG is : 0.00 [g/mol]
Enter the new value, or hit <CR> to leave the same : [<CR>]

Enter data item to be updated : [done]

Current application level data file is LEV
To use an alternate level data file,
enter the file name (<CR> for none)? [<CR>]

NOTE

Editing of the GENERAL data file
has been completed. You are
now entering editing for the
LEVEL data file.

NOTE

An option to let you call in
another LEVEL file to be used
instead of the newly created
file:LEV.DAT. <CR> indicates that
you want to bypass this option.

***** AVAILABLE APPLICATION STUDY REGIONS *****

1. BOSTON MA

CANTON SANDY LOAM FINE S

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

Only one site is available ** automatically selected

***** MODIFICATION OF ANNUAL APPLICATION PARAMETERS *****

1. AR [cm2] surface area
2. Z [m] depth to groundwater table
3. DU [cm] depth of upper unsaturated soil zone
4. DM [cm] depth of middle unsaturated soil zone
5. FRN Freundlich equation exponent
6. PH pH of upper unsaturated soil zone
7. A2PH pH of middle unsaturated soil zone
8. APH ratio: pH lower/upper unsaturated soil zone
9. A2KDE ratio: biodes rate lower/middle soil zone
10. AKDE ratio: biodes rate lower/upper soil zone
11. A2OC ratio: organic carbon lower/middle
12. AOC ratio: organic carbon lower/middle
13. A2CC ratio: clay content lower/middle
14. ACC ratio: clay content lower/upper
15. A2CEC ratio: lower/middle soil zone cation exchange capacity
16. ACEC ratio: lower/upper soil zone cation exchange capacity
17. NAME

Enter data item to be updated : [review]

You may review data for the current year or for all years available.

Enter [C]urrent or [A]ll? [all]

1 BOSTON MA	CANTON SANDY LOAM FINE S 1											
- AR,Z,DU,DM,FRN	10000.	2.74	15.24	68.58	1.00							
- PH,A2PH,APH	4.80	1.00	1.00									
- A2KDE,AKDE,A2OC,AOC,A2CC,ACC	1.00	1.00	1.00	1.00	1.00	1.00						
- A2CEC,ACEC	1.00	1.00										
CUM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
CMH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
CLM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
POLINU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
POLINH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
POLINL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
ISRM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
ASL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
TRANSU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
TRANSM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
TRANSL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
SINKU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
SINKH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
SINKL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
LIGCU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
LIGCH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

LIGCL 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00

Press return to continue? [CR]

Enter data item to be updated : [done]

***** MODIFICATION OF MONTHLY APPLICATION PARAMETERS *****

Current YEAR : 1

1. CUM [us/mL] conc pollutant in soil moisture upper zone
2. CMH [us/mL] conc pollutant in soil moisture middle zone
3. CLM [us/mL] conc pollutant in soil moisture lower zone
4. POLINU [us/cm2] pollution load per unit area entering upper zone
5. POLINH [us/cm2] pollution load per unit area entering middle zone
6. POLINL [us/cm2] pollution load per unit area entering lower zone
7. ISRM [us/cm2] index for pollutant appearance in surface runoff
8. ASL ratio: conc of pollutant/maximum solubility in H2O
9. TRANSU [us/cm2] pollutant transformed in upper zone
10. TRANSM [us/cm2] pollutant transformed in middle zone
11. TRANSL [us/cm2] pollutant transformed in lower zone
12. SINKU [us/cm2] pollutant lost by processes not described (upper)
13. SINKH [us/cm2] pollutant lost by processes not described (middle)
14. SINKL [us/cm2] pollutant lost by processes not described (lower)
15. LIGU [us/cm2] ligand mass input to upper zone
16. LIGH [us/cm2] ligand mass input to middle zone
17. LIGL [us/cm2] ligand mass input to lower zone

Enter data item to be updated : [4]

The current values for POLINU [us/cm2] are

1 Oct	2 Nov	3 Dec	4 Jan	5 Feb	6 Mar
0.00	0.00	0.00	0.00	0.00	0.00
7 Apr	8 May	9 June	10 July	11 Aug	12 Sept
0.00	0.00	0.00	0.00	0.00	0.00

Enter the month(s) to alter? [all]

NOTE

 'all' allows you to change all
 months with one constant value.

Enter the new value(s)? [5000.0]

Current YEAR : 1

Enter data item to be updated : [back]

NOTE

'back' allows you to go back to
the the previous parameter group
for re-editing.

***** MODIFICATION OF ANNUAL APPLICATION PARAMETERS *****

1.	AR	[cm2]	surface area
2.	Z	[m]	depth to groundwater table
3.	DU	[cm]	depth of upper unsaturated soil zone
4.	DM	[cm]	depth of middle unsaturated soil zone
5.	FRN		Freundlich equation exponent
6.	PH		pH of upper unsaturated soil zone
7.	A2PH		pH of middle unsaturated soil zone
8.	APH		ratio: pH lower/upper unsaturated soil zone
9.	A2KDE		ratio: biodes rate lower/middle soil zone
10.	AKDE		ratio: biodes rate lower/upper soil zone
11.	A2OC		ratio: organic carbon lower/middle
12.	AOC		ratio: organic carbon lower/middleupper
13.	A2CC		ratio: clay content lower/middle
14.	ACC		ratio: clay content lower/upper
15.	A2CEC		ratio: lower/middle soil zone cation exchange capacity
16.	ACEC		ratio: lower/upper soil zone cation exchange capacity
17.	NAME		

Enter data item to be updated : [1]

The current value for AR is : 10000. [cm2]

Enter the new value, or hit <CR> to leave the same : [1000.]

Enter data item to be updated : [brief]

NOTE

'brief' suppresses the displaying
of the menu parameter descrip-
tions. This will save editing
time for experienced users.

Enter data item to be updated : [done]

***** MODIFICATION OF MONTHLY APPLICATION PARAMETERS *****

Current YEAR : 1

1.	CUM
2.	CMM
3.	CLM

APP.1:SESOIL SAMPLE RUN FOR 1 YEAR USING BOTH SESIN AND SESOILFM(CONT'D)

- 4. POLINU
- 5. POLINH
- 6. POLINL
- 7. ISRM
- 8. ASL
- 9. TRANSU
- 10. TRANSM
- 11. TRANSL
- 12. SINKU
- 13. SINKM
- 14. SINKL
- 15. LIGU
- 16. LIGM
- 17. LIGL

Enter data item to be updated : [done]

NOTE

Since this data group has already
been edited, no additional
changes are made. (Notice that
the parameter descriptions are
suppressed.)

Do you wish to enter data for additional years? [no]

NOTE

A sample run for multiple years
of data is given in Appendix 2.

Would you like to change any exec file parameters? [no]

NOTE

Unless you are experienced in
the use of the EXEC file para-
meters, changes by you are not
recommended, except for making
multiple-year runs as shown in
Appendix 2.

Would you like to enter data for another run? [no]

NOTE

You are encouraged to try this
multiple run option.

Current GENERAL new input default FILENAME is : GENERAL.DAT
Enter your new input FILENAME? [newgen]

NOTE

You can assign a specific name
to the just edited SESOIL-ready
GENERAL file. Here the name
NEWGEN.DAT was assigned. A <CR>
would have assigned the default
name: GENERAL.DAT.

Current Level 3 new input default filename is : LEVEL3.DAT
Enter the name for your new input file? [newlev]

NOTE

You can assign a specific name
to the just edited SESOIL-ready
LEVEL file. Here, the name
NEWLEV.DAT was assigned. A <CR>
would have assigned the default
name LEVEL3.DAT.

Do you want to run SESOIL: [yes]

NOTE

You have the option of running
SESOIL now (by responding 'yes')
or saving the newly edited data-
sets to be run at a later time
(a 'no' response).

Enter BATCH or INTERACTIVE : [i]

NOTE

You have the option to submit
a batch(b) or interactive(i) job.

Enter Y to output sesoil results to GEMS > [y]

NOTE

This is to indicate if the out-
put data management (SESOILFH)
support is requested or not.

Enter GEMS output dataset name > [pcb]

NOTE

If SESOILFM support was requested
you need to give a name(1 to 12
characters) under which the output
dataset is to be stored in GEMS.

Enter tas field > [sesoil run for pcb at boston]

NOTE

This is to give more descriptive
dataset identification. Press
<CR> to bypass.

DSNAME=TESTPCB1
FMCLOS: IOPT= 1
FORTRAN STOP

NOTE

At this point, SESOIL has been
successfully executed using the
two input files: NEWGEN.DAT and
NEWLEV.DAT. The output file is
named SESOUT.DAT.

NOTE

At this point, SESOILFM has
stored selected SESOIL output
parameters into GEMS under the
name: PCB. The link between
SESOIL and AT123D starts below.

ENTER Y TO SEE SESOIL OUTPUT USABLE FOR AT123D > [y]

NOTE

'y' will allow you to review all
the AT123D-related SESOIL output
parameters prior to building up
an AT123D input file. Those
assumptions which require special
consideration are displayed in
brackets.

THE FOLLOWING GIVES SELECTED SESOIL OUTPUTS AND
THEIR CORRESPONDING AT123D INPUT VALUES
(SEE USER'S GUIDE FOR ASSUMPTIONS MADE)

(1)SESOIL-AREA(AR): 1000.000 SQ.CH.
AT123D-SOURCE SIZE: RL1= 0.0000000E+00 RL2= 0.3162278
(METERS) RB1= -0.1581139 RB2= 0.1581139
RH1= 0.0000000E+00 RH2= 0.0000000E+00

(2)SESOIL-EFFECTIVE POROSITY(N): 0.2600000
AT123D-EFFECTIVE POROSITY(POR): 0.2600000

(3)SESOIL-INTR. PERM.(K1): 5.5440285E-08 SQ. CH.
AT123D-HYD. COND.(HCOND): 0.1710666 M/HR
**[NOTE: POROSITY AND HYDR. COND. IN
GROUNDWATER LAYERS MAY DIFFER SIGNIFICANTLY
FROM THOSE USED IN SESOIL LAYERS.]

(4)SESOIL-AD. COEF.(K): 23.00000 ML/G
AT123D-DIS. COEF.(AKD): 2.3000002E-02 M**3/KG
**[NOTE: K IS DETERMINED INDIRECTLY
FROM KOC AND OC.]
**[NOTE: ACTUAL VALUE OF AKD MAY BE LOWER
DUE TO LOWER OC CONCENTRATION (<0.1%)
IN GROUNDWATER LAYER OF AT123D.]

(5)SESOIL-HYDROL. RATE(KNH): 0.0000000E+00 1/DAY
AT123D-DECAY RATE(RAMADA): 0.0000000E+00 1/HR
**[NOTE: OTHER PROCESSES IN ADDITION TO
NEUTRAL HYDROLYSIS MAY BE CONSIDERED FOR
TOTAL DEGREDDATION IN GROUNDWATER.]

(6)SESOIL-SOIL DENSITY(RS): 1.350000 G/CM**3
AT123D-BULK DENSITY(RHOB): 1350.000 KG/M**3

(7)SESOIL-MASS INTO GRW(12 VALUES PER YEAR)(KG/HR):

1	0.0000E+00	0.0000E+00	0.0000E+00	0.0000E+00	0.2329E-11	0.4197E-11
	0.4553E-11	0.5116E-11	0.5013E-11	0.4576E-11	0.9528E-11	0.1260E-10

ENTER Y TO CREATE AT123D INPUT FILE
USING THE SELECTED SESOIL OUTPUTS > [y]

NOTE

'y' allows you to create an
AT123D input file based on the
above SESOIL simulation run
outputs.

ENTER THE DATA FILE NAME(6 CHARACTERS) > [pcb001]

NOTE

An AT123D input data file name
has to be assigned here.
(6 characters are required.)

THE USABLE SESOIL VALUES HAVE BEEN STORED
IN AN AT123D INPUT DATASET: pcb001.DAT
NOTE: THIS DATASET IS ALSO FILLED WITH OTHER
DEFAULT VALUES WHICH ARE TO BE USED WITH CAUTION.
YOU MAY EDIT THOSE DEFAULT VALUES USING AT123DIN.

DO YOU WANT TO RUN AT123D ? : [no]

NOTE

'yes' will allow you to run
AT123D using AT123DIN and
AT123DOUT. See the AT123D
User's Guide(GSC,1984).

ENTER AN OPERATION:

NOTE

This is back to GEMS.

APPENDIX 2: SESOIL SAMPLE RUN FOR MULTIPLE YEARS USING SESIN

Enter an operation: [m]

Enter a model name: [sesoil]

Are you creating SESOIL input files: [yes]

SESIN
INTERACTIVE SESOIL INPUT
GSC VER1.1
6-DEC-84

For a list of commands type : HELP COMMANDS

Do you want to use data from the cities database? [no]

Current GENERAL data file is : DBAO:[CEEDDAL.DAT]GENERAL.DAT

To use an alternate GENERAL data file,

enter the file name (<CR> for none)? [newsen]

NOTE

Here, the data file:NEWGEN.DAT
created in Appendix 1 is called
in instead of accessing the 70-
cities database. (A <CR> entry
to the last prompt would have
accessed the default GENERAL data
file shown in Appendix 4.

***** DETERMINATION OF SESOIL POLLUTANT APPLICATION LEVEL *****

Are pollutant loadings available as monthly or annual data? [m]

Are there 2 or 3 unsaturated soil layers? [3]

NOTE

The two above prompts determine
the LEVEL type file used.

Selected SESOIL application level is : 3

***** AVAILABLE CLIMATIC REGIONS *****

1. BOSTON MA

CANTON SANDY LOAM FINE S

Only one site is available ** automatically selected

***** CLIMATOLOGICAL PARAMETERS *****

1. L [deg N] latitude of area
2. NAME

Enter data item to be updated : [review]

You may review data for the current year or for all years available.
Enter [C]urrent or [A]ll? [all]

1 BOSTON MA	CANTON SANDY LOAM FINE S												1
L	42.36												
TA	13.00	7.330	5.556	-1.56	-0.89	3.39	9.22	14.78	20.00	24.00	21.83	18.06	
NN	0.53	0.6	0.6	0.6	0.56	0.56	0.58	0.58	0.57	0.56	0.53	0.51	
S	0.57	0.61	0.61	0.57	0.57	0.57	0.53	0.57	0.6	0.56	0.57	0.61	
A	0.18	0.21	0.28	0.37	0.35	0.28	0.21	0.18	0.18	0.18	0.18	0.18	
REP	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	
MPM	7.67	11.46	10.77	9.37	8.99	10.19	8.86	8.81	8.10	6.96	8.79	8.03	
MTR	0.6667	1.040	0.7083	0.375	0.750	0.2917	0.3750	0.3333	0.2083	0.5417	0.0833	0.5	
MN	7.00	4.00	9.00	6.00	8.00	5.00	9.00	7.00	8.00	9.00	5.00	7.00	
MT	9.00	11.00	12.00	12.00	11.00	12.00	11.00	11.00	11.00	9.00	10.00	9.00	

Press return to continue? [CR]

Enter data item to be updated : [brief]

NOTE

'brief' is used for more efficient
editing.

Enter data item to be updated : [done]

Current YEAR : 1

1. TA
2. NN
3. S
4. A
5. REP
6. MPM
7. MTR
8. MN
9. MT

Enter data item to be updated : [done]

APP.2: SESOIL SAMPLE RUN FOR MULTIPLE YEARS USING SESIN(CONT'D)

Would you like to enter data for another year? [yes]
The previous years data is available. Do you wish to use it? [yes]

NOTE

A second year of data for the
GENERAL file is generated by
repeating the first year of data.

Current YEAR : 2

1. TA
2. HN
3. S
4. A
5. REP
6. MPM
7. MTR
8. MN
9. MT

Enter data item to be updated : [done]

NOTE

This second year of GENERAL data
can be edited. Here, no changes
are made.

Would you like to enter data for another year? [no]

NOTE

Additional years of data can be
generated in the same way up to
10 years. Here, only one addi-
tional year has been generated.

***** AVAILABLE SOILS *****

1. CANTON SANDY LOAM FINE SANDY LOAM BOSTON

Only one soil is available ** automatically selected

***** SOIL PARAMETERS *****

APP.2: SESOIL SAMPLE RUN FOR MULTIPLE YEARS USING SESIN(CONT'D)

1. RS
2. K1
3. C
4. N
5. OC
6. CC
7. CEC
8. K1U
9. K1M
10. K1L
11. RDUST
12. NAME

Enter data item to be updated : [done]

***** AVAILABLE CHEMICALS *****

1. PCB-C12H9CL

Only one chemical is available ** automatically selected

***** CHEMICAL PARAMETERS *****

1. SL
2. KOC
3. DA
4. KDE
5. H
6. K
7. MWT
8. VAL
9. KNH
10. KBH
11. KAH
12. SK
13. B
14. MWTLIG
15. NAME

Enter data item to be updated : [review]

1 PCB-C12H9CL					
SL,KOC,DA,KDE,HENRY,K	1.17	11500.	0.06	0.0346	0.0059 0.00
MWT,VAL,KNH,KBH,KAH	133.40	0.00	0.00	0.00	0.00
SK,B,MWTLIG	0.00	0.00	0.00		

Press return to continue? [<CR>]

Enter data item to be updated : [done]

APP.2: SESOIL SAMPLE RUN FOR MULTIPLE YEARS USING SESIN(CONT'D)

NOTE

No changes have been made to any
of the GENERAL data parameters.

Current application level data file is DBA0:[EEDDAL.DAT]LEVEL3.DAT
To use an alternate level data file,
enter the file name (<CR> for none)? [newlev]

NOTE

Here, the data file:NEWLEV.DAT
created in the sample run in
Appendix 1 is called in instead
of accessing the 70-cities data-
base. (A <CR> entry would have
accessed one of the default level
data files shown in Appendix 4.)

***** AVAILABLE APPLICATION STUDY REGIONS *****

1. BOSTON MA CANTON SANDY LOAM FINE S

Only one site is available ** automatically selected

***** MODIFICATION OF ANNUAL APPLICATION PARAMETERS *****

1. AR
2. Z
3. DU
4. DM
5. FRN
6. PH
7. A2PH
8. APH
9. A2KDE
10. AKDE
11. A2OC
12. AOC
13. A2CC
14. ACC
15. A2CEC
16. ACEC
17. NAME

Enter data item to be updated : [done]

APP.2: SESOIL SAMPLE RUN FOR MULTIPLE YEARS USING SESIN(CONT'D)

***** MODIFICATION OF MONTHLY APPLICATION PARAMETERS *****

Current YEAR : 1

1. CUM
2. CHM
3. CLM
4. POLINU
5. POLINH
6. POLINL
7. ISRM
8. ASL
9. TRANSU
10. TRANSH
11. TRANSL
12. SINKU
13. SINKH
14. SINKL
15. LIGU
16. LIGH
17. LIGL

Enter data item to be updated : [review]

You may review data for the current year or for all years available.
Enter [C]urrent or [A]ll? [all]

1 BOSTON MA	CANTON SANDY LOAM FINE S 1											
- AR,Z,DU,DM,FRN	10000.	2.74	15.24	68.58	1.00							
- PH,A2PH,APH	4.80	1.00	1.00									
- A2KDE,AKDE,A2DC,AOC,A2CC,ACC	1.00	1.00	1.00	1.00	1.00	1.00						
- A2CEC,ACEC	1.00	1.00										
CUM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
CHM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
CLM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
POLINU	5000.	5000.	5000.	5000.	5000.	5000.	5000.	5000.	5000.	5000.	5000.	5000.
POLINH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
POLINL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
ISRM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
ASL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
TRANSU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
TRANSH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
TRANSL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
SINKU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
SINKH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
SINKL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
LIGU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
LIGH	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
LIGL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00

APP.2: SESOIL SAMPLE RUN FOR MULTIPLE YEARS USING SESIN(CONT'D)

Press return to continue? [CR]

Current YEAR : 1

Enter data item to be updated : [done]

Do you wish to enter data for additional years? [yes]

The previous years data is available. Do you wish to use it? [yes]

NOTE

A second year of data for the
LEVEL data file is generated by
repeating the first year of data.

***** MODIFICATION OF MONTHLY APPLICATION PARAMETERS *****

Current YEAR : 2

1. CUM
2. CMH
3. CLM
4. POLINU
5. POLINM
6. POLINL
7. ISRM
8. ASL
9. TRANSU
10. TRANSM
11. TRANSL
12. SINKU
13. SINKM
14. SINKL
15. LIGU
16. LIGM
17. LIGL

Enter data item to be updated : [nobrief]

NOTE

'nobrief' turns off the 'brief'
command and all parameter descri-
ptions will be displayed.

Current YEAR : 2

Enter data item to be updated : [done]

NOTE

No changes have been made to the
generated second year of LEVEL
data.

Do you wish to enter data for additional years? [no]

NOTE

Note that the number of years of
data in the GENERAL and LEVEL
data files must be the same.

Would you like to change any exec file parameters? [yes]

NOTE

A 'yes' response here displays
the EXEC file menu for editing.

1. JRUN Do Not Change!
2. LEVEL level of operation
3. JRE index of area of application (GENERAL file)
4. JSO soil index
5. JCH chemical index
6. JNUT nutrient cycle (cannot change)
7. JAPFL application area (LEVEL file)
8. JYRS number of years to be simulated (must be 1 for annual)

Enter data item to be updated : [8]

The current value for JYRS is : 2.00

Enter the new value, or hit <CR> to leave the same : [5]

NOTE

Here, it is desired to increase
the number of years to be simulated
from 2 to 5. Because only 2 years
of data are available (as previously
created), the auxiliary pollution
load data file is created as follows.
(This is the decision to enter
Branch 1-A! See Fig. 1)

NOTE

This method of generating additional data is usually used for long periods of simulation. Up to 100 years may be simulated. Note that only the pollution load parameters (POLINU, POLINM, and POLINL) are edited. The remaining GENERAL and LEVEL data parameters are automatically generated based on repeated use of the last year of available data.

Enter data item to be updated : [done]

AUX LOADINGS FILE

You have 2 years of data.
you must enter 3 more years.

- | | |
|------------|--|
| 1. SAME | sets all elements to the same number |
| 2. RECYCLE | recycles existing loading data for remaining years |
| 3. EDIT | normal SESIN data entry/editing procedures |

NOTE

Two options are available for generating the auxiliary loading data file. (Option no. 2 has not yet been implemented in the present version.)

Enter method of loading entry: [1]

NOTE

The 'SAME' method of generating loading data repeats a 1-year loading scenario for all the additional years of data required.

Enter soil layer (1, 2, or 3)? [1]

NOTE

Here, the loading is applied to the upper soil layer (layer 1), (POLINU).

Enter month(s) to alter? [all]

NOTE

Here, all months in a year will
have the same loading value.

Enter loadings? [2000.]

NOTE

Here, a loading value of 2000 is
applied to all months of all of
the remaining years required
(years 3,4 and 5).

You have 5 years of data,
you must enter 0 more years.

- | | |
|------------|--|
| 1. SAME | sets all elements to the same number |
| 2. RECYCLE | recycles existing loading data for remaining years |
| 3. EDIT | normal SESIN data entry/editing procedures |

Enter method of loading entry: [3]

NOTE

The 'EDIT' method of loading data
entry allows you to edit the
loading quantities for individual
years.

***** ENTRY OF AUXILIARY LOADING PARAMETERS *****

Current YEAR : 3

- | | |
|-----------|-----------------------|
| 1. POLINU | Upper level loadings |
| 2. POLINM | middle level loadings |
| 3. POLINL | lower level loadings |

NOTE

Editing begins with the first
year in the AUX loading file
(year 3).

APP.2: SESOIL SAMPLE RUN FOR MULTIPLE YEARS USING SESIN(CONT'D)

Enter data item to be updated : [1]

NOTE

 Here, the source load is applied
 to the upper soil layer (POLINU).

The current values for POLINU are

1 Oct	2 Nov	3 Dec	4 Jan	5 Feb	6 Mar
2000.00	2000.00	2000.00	2000.00	2000.00	2000.00
7 Apr	8 May	9 June	10 July	11 Aug	12 Sept
2000.00	2000.00	2000.00	2000.00	2000.00	2000.00

Enter the month(s) to alter? [done]

NOTE

 No months are to be altered.

Current YEAR : 3

Enter data item to be updated : [next]

NOTE

 'next' advances the editing to
 the next year.

Current YEAR : 4

Enter data item to be updated : [1]

The current values for POLINU are

1 Oct	2 Nov	3 Dec	4 Jan	5 Feb	6 Mar
2000.00	2000.00	2000.00	2000.00	2000.00	2000.00
7 Apr	8 May	9 June	10 July	11 Aug	12 Sept
2000.00	2000.00	2000.00	2000.00	2000.00	2000.00

Enter the month(s) to alter? [all]

Enter the new value(s)? [1000.]

NOTE

 Here, the values of POLINU for
 all months of year 4 are changed

from 2000 to 1000.

Current YEAR : 4

Enter data item to be updated : [done]

NOTE

No more changes are to be made
for year 4.

You have 5 years of data.
you must enter 0 more years.

1. SAME sets all elements to the same number
2. RECYCLE recycles existing loading data for remaining years
3. EDIT normal SESIN data entry/editing procedures

Enter method of loading entry: [done]

NOTE

Here, 'done' indicates that you
have completed all editing of
the auxiliary loading data file
AUX.

Would you like to enter data for another run? [no]

Current GENERAL new input default FILENAME is : GENERAL.DAT
Enter your new input FILENAME? [newsen2]

Current Level 3 new input default filename is : LEVEL3.DAT
Enter the name for your new input file? [newlev2]
Enter filename for aux loadings file: [aux5]

NOTE

You must also specify a filename
for the auxiliary loadings file.

Do you want to run SESOIL: [yes]

Enter BATCH or INTERACTIVE : [b]

Enter Y to output sesoil results to GEMS > [<CR>]
Job 9410 entered on queue SYS BATCH

NOTE

Here, the run is submitted as a
batch job.

Enter an operation:

NOTE

This is back to GEMS.

APPENDIX 3: SESOIL SAMPLE RUN BYPASSING BOTH SESIN AND SESOILFM

Enter an operation: [m]

Enter a model name: [sesoil]

Are you creating SESOIL input files: [no]

NOTE

This is the decision to enter
either Branch 1 or Branch 2(See
Fig. 1). This example shows
that Branch 2(i.e. no SESIN
support) has been requested.

Enter DUMMY below if the file(s) is not needed
by your run

Enter filename of GENERAL data file: [newgen]

Enter filename of LEVEL0 data file: [dummy]

Enter filename of LEVEL1 data file: [dummy]

Enter filename of LEVEL2 data file: [dummy]

Enter filename of LEVEL3 data file: [newlev]

Enter filename of EXEC data file: [exec]

Enter filename of AUXILIARY data file(needed if JYRS>run years): [dummy]

NOTE

Since SESIN support is not
requested, the above are to input
your own GENERAL, LEVEL, and EXEC
data files. Only one LEVEL file
is needed for each run. The
AUXILIARY data file is needed if
the value of JYRS in the EXEC
file is greater than the number
of years of data available in
the GENERAL and LEVEL files
(see Appendix 2).

Enter BATCH or INTERACTIVE : [i]

Enter Y to output SESOIL results to GEMS > [<CR>]

NOTE

<CR> indicates that SESOILFM
support is not requested.

DSNAME=-1

FMCLIB: IOPT=

0

FORTRAN STOP

NOTE

Indication of SESOIL FORTRAN
execution completion.

ENTER Y TO SEE SESOIL OUTPUT USABLE FOR AT123D > [<CR>]

NOTE

Review of SESOIL output usable
for AT123D is not requested.

ENTER Y TO CREATE AT123D INPUT FILE
USING THE SELECTED SESOIL OUTPUTS > [<CR>]

NOTE

Build up of AT123D input file
via SESOIL output is not requested.

DO YOU WANT TO RUN AT123D ?; [no]

The output file for this modeling run is named SESOUT.DAT
Enter rename, print, type or <CR> to exit: [<CR>]

NOTE

Here, the decision is to exit.

Enter a model name:

NOTE

This is back to GEMS.

*****NOTE*****

In summary, this example shows a SESOIL run without the support
of SESIN and SESOILFM. The user needs to input his own data
files. The complete SESOIL output is stored in a file named
SESOUT.DAT.

APPENDIX 4: SESIN-PROVIDED DEFAULT DATA FILES: GENERAL, LEVEL0, LEVEL1,
LEVEL2, AND LEVEL3

GENERAL DATA FILE

1 REGIONAL DESCRIPTIONS; CLIMATIC-STORM DATA: (TEST DATA ONLY)

1 CLINTON, MA. (SUB-HUMID):

- L,TA,NN,S,A,REP	42.50	8.40	0.35	0.70	0.30	0.00
- MPA,MTR,MN,HT	94.10	0.32	109.00	365.00		
- MP(M) (OCTOBER-MARCH)	10.36	8.96	12.52	11.75	8.33	8.33
- MP(M) (APRIL-SEPTEMBER)	8.38	7.82	6.60	5.08	6.04	8.46

2 SITE B (MONTANA) (1940-1979)

- L,TA,NN,S,A,REP	45.00	10.05	0.30	.55	0.25	0.00
- MPA,MTR,MN,MT	36.74	0.24	103.0	365.00		
- MP(M)OCT-MAR	2.00	1.63	2.59	4.11	5.44	6.25
- MP(M)APR-SEP	2.33	2.36	3.40	2.84	1.96	1.83

3 SITE A (KANSAS) (1940-1990)

- L,TA,NN,S,A,REP	39.00	14.30	0.30	60.0	0.10	0.00
- MPA,MTR,MN,MT	84.83	0.20	105.0	365.00		
- MP(M)OCT-MAR	6.40	4.22	2.95	2.46	3.20	5.20
- MP(M)APR-SEP	7.82	11.10	12.17	9.97	10.24	9.12

4 CLINTON,MASS.(40 YR AVERAGED DATA FOR TEST ONLY) 10

L	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50
TA	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40
NN	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
S	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70
A	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3
REP	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
MPM	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84
MTR	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
MN	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08
MT	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42

L-2	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50
TA-2	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40
NN-2	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
S-2	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70
A-2	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3
REP-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
MPM-2	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84
MTR-2	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
MN-2	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08
MT-2	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42

L-3	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50
TA-3	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40
NN-3	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
S-3	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70
A-3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3
REP-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
MPM-3	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84

APP. 4: SESIN-PROVIDED DEFAULT DATA FILES (CONT'D)

[illegible]

APP. 4: SESIN-PROVIDED DEFAULT DATA FILES (CONT'D)

REP-8	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
MPM-8	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84
MTR-8	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
MN-8	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08
MT-8	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42
L-9	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50
TA-9	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40
MN-9	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
S-9	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70
A-9	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3
REP-9	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
MPM-9	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84
MTR-9	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
MN-9	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08
MT-9	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42
L-10	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50	42.50
TA-10	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40	8.40
MN-10	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
S-10	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70	.70
A-10	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3
REP-10	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
MPM-10	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84	7.84
MTR-10	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32	.32
MN-10	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08	9.08
MT-10	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42	30.42

2 SOIL CLASSIFICATION DATA: (TEST DATA ONLY)

1 CLAY

- RS,K1,C,N,OC,CC	1.32	1.E-10	12.00	0.45	1.46
- CEC,K1U,K1M,K1L	15.00	0.00	0.00	0.00	

2 SILTY-LOAM-KANSAS

- RS,K1,C,N,OC,CC	1.32	7.E-09	6.00	0.35	2.9
- CEC	0.00				

3 SILTY-CLAY-MONTANA

- RS,K1,C,N,OC,CC	1.324.2E-10	10.00	0.35	.63	
- CEC	0.00				

4 SILTY-LOAM-KANSAS (K1 ADJUSTED)

- RS,K1,C,N,OC,CC	1.32	1.E-13	6.00	0.35	2.9
- CEC	0.00				

5 CLAY-LOAM GENERIC

-RS,K1,C,N,OC,CC	1.32	.29E-9	10.00	0.35	1.3 .20E-1
-CEC	.00				

6 SANDY-LOAM

- RS,K1,C,N,OC,CC	1.322.5E-10	4.00	0.25	0.10	0.00
- CEC	0.00				

7 TEST SOIL

- RS,K1,C,N,OC,CC	1.32	.25E-9	4.00	0.25	0.10	5.00
- CEC,K1U,K1M,K1L	15.00	.22E-9	.33E-9	.44E-9		

8 SILTY-LOAM GENERIC

APP. 4: SESIN-PROVIDED DEFAULT DATA FILES (CONT'D)

	-RS,K1,C,N,OC,CC	1.32	1.2E-9	6.00	0.35	2.9
	-CEC	0.00				
9	SAND-GENERIC					
	-RS,K1,C,N,OC,CC	1.32	1.0E-7	3.50	0.25	.01
	-CEC	0.00				
3	CHEMISTRY DATA; (TEST DATA ONLY)	<i>Please review & modify before use!</i>				
1	COPPER					
	SL,KOC,DA,KDE,H,K	0.00	0.00	0.00	0.00	10.00
	MWT,VAL,KNH,KBH,KAH	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
2	SODIUM					
	SL,KOC,DA,KDE,H,K	0.00	0.00	0.00	0.00	.10
	MWT,VAL,KNH,KBH,KAH	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
3	CHROMIUM					
	SL,KOC,DA,KDE,H,K	0.00	0.00	0.00	0.00	250.00
	MWT,VAL,KNH,KBH,KAH	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
4	NICKEL					
	SL,KOC,DA,KDE,H,K	0.00	0.00	0.00	0.00	2.50
	MWT,VAL,KNH,KBH,KAH	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
5	NAPHTHALENE					
	SL,KOC,DA,KDE,H,K	0.0	1070.006.98E-6	0.00	0.00	0.00
	MWT,VAL,KNH,KBH,KAH	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
6	ANTHRACENE					
	SL,KOC,DA,KDE,H,K	0.00	1.6E45.89E-6	0.00	0.00	.00
	MWT,VAL,KNH,KBH,KAH	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
7	SODIUM-ADJUSTED KOC					
	SL,KOC,DA,KDE,H,K	0.00	0.00	0.00	0.00	1.00
	MWT,VAL,KNH,KBH,KAH	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
8	PERCHLOROETHYLENE(=TETRACHLOROETHYLENE)					
	SL,KOC,KD,KDE,HENRY,K	150.00	468.00	.0746	0.00	.00204
	MWT,VAL,KNH,KBH,KAH,MWTLIG	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
9	1,1,1-TRICHLOROETHANE					
	SL,KOC,DA,KDE,HENRY,K	720.00	174.00	.0803	0.00	.00298
	MWT,VAL,KNH,KBH,KAH,MWTLIG	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
10	METHYLCHLOROFORM					
	SL,KOC,DA,KDE,HENRY,K	740.00	189.00	.0825	0.00	.00231
	MWT,VAL,KNH,KBH,KAH,MWTLIG	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
11	METHYLENE CHLORIDE					
	SL,KOC,DA,KDE,HENRY,K	13200.	11.00	.1033	0.003.68E-4	0.00
	MWT,VAL,KNH,KBH,KAH,MWTLIG	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00
12	CARBON TETRACHLORIDE					
	SL,KOC,DA,KDE,HENRY,K	785.00	269.00	.0828	0.00	.00232
	MWT,VAL,KNH,KBH,KAH,MWTLIG	0.00	0.00	0.00	0.00	0.00

APP. 4: SESIN-PROVIDED DEFAULT DATA FILES (CONT'D)

	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00	0.00
13	FREON 113						
	SL,KOC,DA,KDE,HENRY,K	170.00	642.00	.0693	0.00	.391	0.00
	MWT,VAL,KNH,KBH,KAH,MWTLIG	0.00	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00	0.00
14	TRICHLOROETHYLENE (TCE)						
	SL,KOC,DA,KDE,HENRY,K	1100.00	100.00	.0830	0.009.44E-3		0.00
	MWT,VAL,KNH,KBH,KAH,MWTLIG	0.00	0.00	0.00	0.00	0.00	0.00
	SK,B,MWTLIG	0.00	0.00	0.00	0.00	0.00	0.00
15	TEST CHEM						
	SL,KOC,DA,KDE,HENRY,K	1100.00	360.00	8.3E-6	0.2	9.44E-3	
	MWT,VAL,KNH,KBH,KAH	5.00	2.00	.7E-5	8.00	9.00	10.00
	SK,B,MWTLIG	100000.	2.00	10.00	0.00	0.00	0.00
9	END FILE						

LEVEL0 DATA FILE

1	CLINTON, MA. (SUB-HUMID): FLAG						
	- AR,Z,DU,PH,APH	1.00	100.0	15.00	7.00	1.00	
	- AKDE,AOC,ACC,ISRA,ASL,ACEC	.50	.50	0.50	1.00	.001	
	- POLINU,POLINL,LIGU,LIGL	200.0	1000.0	1000.0	2000.0		
	- THA,INF,RGA,RSA	25.7	60.70	26.2	33.9		
2	TEST LOCATION						
	- AR,Z,DU,PH,APH	1.00	100.00	15.00	8.00	.875	
	- AKDE,AOC,ACC,ISRA,ASL,ACEC	1.10	1.20	1.30	.40	0.5	.001
	- POLINU,POLINL,LIGC,ALIGC	200.00	1000.0	10.00	50.00		
	- THA,IA,RGA,RSA	25.70	60.70	26.20	33.90		
9	END FILE						

LEVEL1 DATA FILE

1	CLINTON, MA. (SUB-HUMID):						
	- AR,Z,DU,PH,APH	1.00	100.0	15.00	7.00	1.00	
	- AKDE,AOC,ACC,ISRA,ASL,ACEC	.50	.50	0.50	1.00	.001	
	- SPOLI,SPOLL,LIGU,LIGL	200.0	1000.0	1000.0	2000.0		
2	TEST LOCATION						
	- AR,Z,DU,PH,APH	10.00	100.00	15.00	8.00	.875	
	- AKDE,AOC,ACC,ISRA,ASL,ACEC	1.10	1.20	1.30	.40	0.5	.001
	- POLINU,POLINL,LIGC,LIGL	200.00	1000.0	10.00	50.00		
9	END OF FILE						

LEVEL2 DATA FILE

1 KANSAS-COPPER (OCT. '78-SEP '80)						2						
- AR,Z,DU,PH,APH						1.00	50.0	15.00				
- AKDE,ADC,ACC,FRN,ACEC						0.00	1.00	0.00	1.00	0.00		
CUM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CLM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINU	0.00	0.00	0.00	0.00	0.00	0.00	3.24	0.00	0.00	0.00	0.00	0.0
POLINL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ISRM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ASL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CUM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CLM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ISRM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ASL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
2 TEST RUN (CLINTON, MA. 40 YEARS)						5						
- AR,Z,DU,PH,APH						1.00	100.00	15.00	8.00	.975		
- AKDE,ADC,ACC,FRN,ACEC						1.10	1.20	1.30	1.40	.001		
CUM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CLM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINU	200.0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINL	1000.0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ISRM	1.00	1.00	1.00	1.00	1.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSU	0.20	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSL	0.30	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKU	0.40	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKL	0.50	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGU	1000.0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGL	2000.0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CUM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CLM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ISRM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0

APP. 4: SESIN-PROVIDED DEFAULT DATA FILES (CONT'D)

9 END FILE

LEVEL3 DATA FILE

1 TEST RUN						10						
-AR,Z,DU,DM,FRN						1.00	100.00	15.00	0.01	1.40		
-PH,A2PH,APH						8.00	1.00	.875				
-A2KDE,AKDE,A2OC,AOC,A2CC,ACC						1.10	1.10	1.20	1.20	1.30	1.30	
-A2CEC,ACEC						1.30	1.30					
CUM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CHM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CLM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINU	200.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINL	1000.0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ISRM	1.00	1.00	1.00	1.00	1.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ASL	.001	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSU	0.20	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSL	0.30	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKU	0.40	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKL	0.50	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGCU	1000.0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGCM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGCL	2000.0	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CUM-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CHM-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CLM-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ISRM-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ASL-2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSU	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSM	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
TRANSL	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
SINKL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGCU2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGCM2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
LIGCL2	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CUM-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CHM-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
CLM-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINU3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINM3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
POLINL3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ISRM-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0
ASL-3	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.0

APP. 4: SESIN-PROVIDED DEFAULT DATA FILES (CONT'D)

[illegible]

APP. 4: SESIN-PROVIDED DEFAULT DATA FILES (CONT'D)

[illegible]